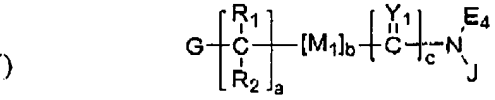
Amendments to the Claims:

This listing of claims will replace all prior versions and listings of claims in the application:

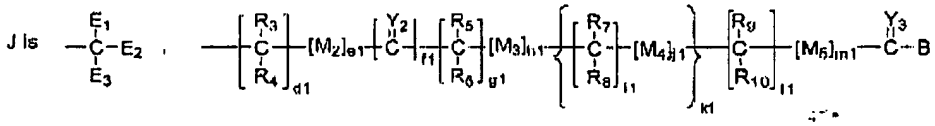
Listing of Claims:

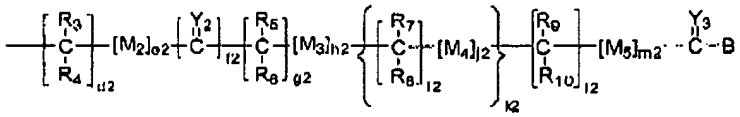
1. (currently amended) A compound comprising the formula:



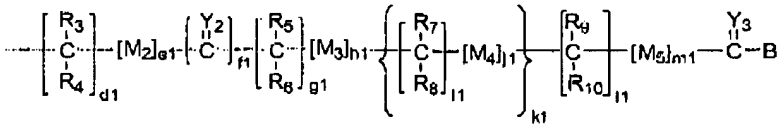
(I)

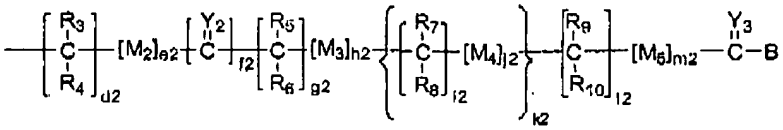
wherein:

J is 

and or 

E_{1-4} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} heteroalkyls, C_{1-6} alkoxy, phenoxy, C_{1-6} heteroalkoxy,

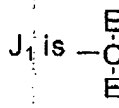


and or 

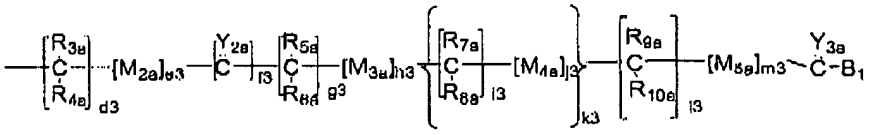
and at least one of E_{1-4} includes a B moiety;

B is a leaving group, OH, a residue of a hydroxyl-containing moiety, a residue of an amine-containing moiety or

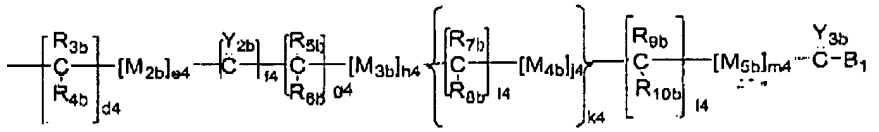
wherein E_5 is independently selected from the same group which defines E_{1-4};

J_1 is  ,

E_{1a-3a} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} heteroalkyls, C_{1-6} alkoxy, phenoxy, C_{1-6} heteroalkoxy,

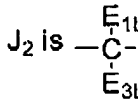


and or

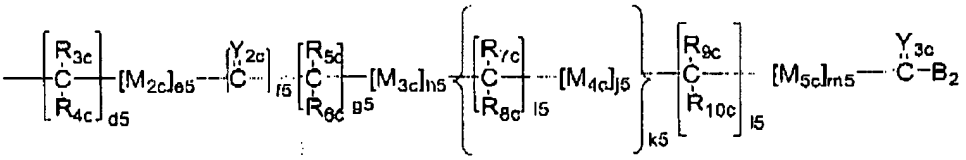


wherein B_1 is a leaving group, OH, a residue of a hydroxyl-containing moiety or a residue of an amine-containing moiety or

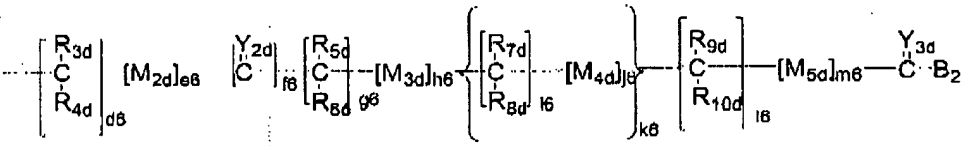
wherein E_6 is independently selected from the same group which defines E_{1-4};

J_2 is  ,

wherein E_{1b-3b} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} heteroalkyls, C_{1-6} alkoxy, phenoxy, C_{1-6} heteroalkoxy,



or



wherein B_2 is a leaving group, OH, a residue of a hydroxyl-containing moiety or a residue of an amine-containing moiety;

G is a polymeric residue;

Y_{1-3}, Y_{2a-d} and Y_{3a-d} are each independently O, S or NR_{11a}

M_{1-4}, M_{2a-2d}, M_{3a-3d}, and M_{4a-4d} are each independently O, S or NR_{11b};

M_5 and M_{5a-d} are each independently X or Q,

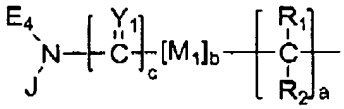
wherein X is an electron withdrawing group and Q is a moiety containing a free electron pair positioned three to six atoms from $C(=Y_3)$ or $C(=Y_{3a-d})$;

R_{1-10}, R_{1a-11a}, R_{1b-11b}, R_{1c-10c} and R_{1d-10d} are each independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} heteroalkyls, C_{1-6} alkoxy, phenoxy and C_{1-6} heteroalkoxy; ~~and~~

a, b, c, d1- d6, e1-e6, f1- f6, g1- g6, h1- h6, ~~i1- i6, j1- j6, k1- k6,~~ l1- l6, m1- m6 are each independently zero or a positive integer; and

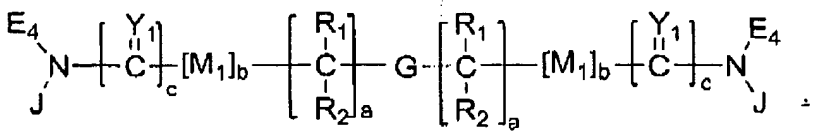
i1-i6, j1-j6 and *k1-k6* are each independently selected positive integers.

2. (original) The compound of claim 1, wherein G further comprises a capping group A, which is selected from the group consisting of hydrogen, CO_2H, C_{1-6} alkyl moieties, and

(1') 

wherein *a*, *b*, *c*, R_{1-2}, M_1, Y_1, E_4 and J are the same as set forth in claim 1.

3. (currently amended) A compound of claim 2, of the formula:



4. (currently amended) The compound of claim 1, where *a, b, c, d1- d6, e1-e6, f1- f6, g1- g6, h1- h6, ~~i1- i6, j1- j6, k1- k6,~~ l1- l6,* and *m1- m6* are independently zero, one or two; and *i1- i6, j1- j6,* and *k1-k6* are independently one or two.

5. (original) The compound of claim 1, wherein R_1 and R_2 are both H, *a* and *c* are one, Y_1 is O and both E_1 and E_4 are H.

6. (original) The compound of claim 1, wherein G is polyalkylene oxide residue.

7. (original) The compound of claim 6, wherein G is a polyethylene glycol residue.

8. (original) The compound of claim 1, wherein G is $-O-(CH_2CH_2O)_x$ or $-O-(CH(CH_3)CH_2O)_x$,

wherein x is the degree of polymerization.

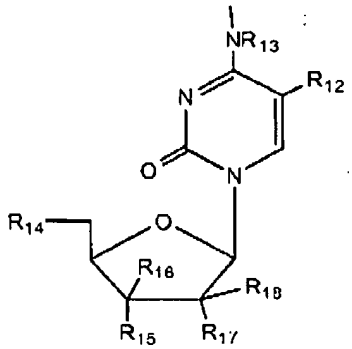
9. (currently amended) The compound of claim 8, wherein G is $-O-(CH_2CH_2O)_x$ and x is a positive integer so that the weight average molecular weight is at least about 20,000 daltons.

10. (currently amended) The compound of claim 9, wherein G has a weight average molecular weight of from about 20,000 to about 100,000 daltons.

11. (currently amended) The compound of claim 10, wherein G has a weight average molecular weight of from about 25,000 to about 60,000 daltons.

12. (original)The compound of claim 1, wherein B is a residue of an amine -containing moiety.

13. (original) The compound of claim 12, wherein said amine-containing moiety is

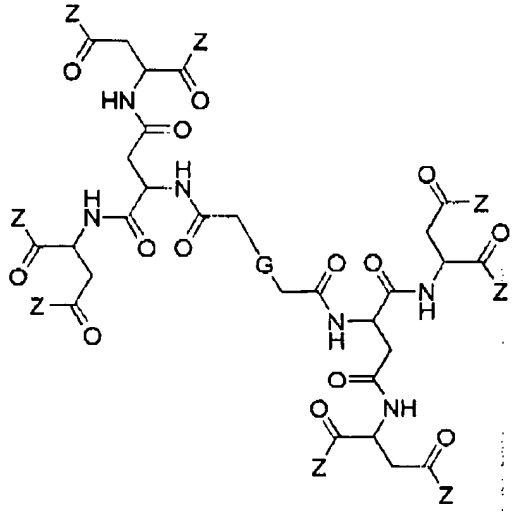


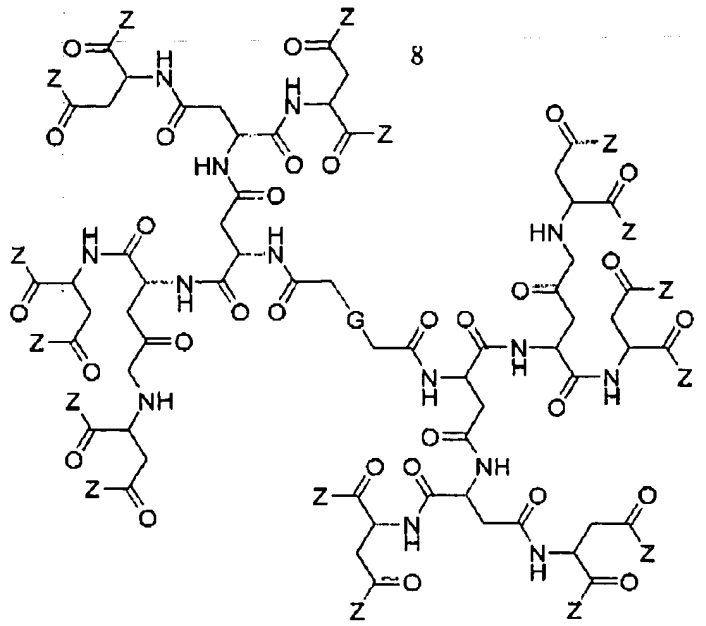
wherein

$R_{12\text{-}13}$ are independently selected from the group consisting of hydrogen, $C_{1\text{-}6}$ alkyls, $C_{3\text{-}12}$ branched alkyls, $C_{3\text{-}8}$ cycloalkyls, $C_{1\text{-}6}$ substituted alkyls, $C_{3\text{-}8}$ substituted cycloalkyls, aryls, halo, substituted aryls, aralkyls, $C_{1\text{-}6}$ heteroalkyls, substituted $C_{1\text{-}6}$ heteroalkyls;

$R_{14\text{-}18}$ are independently selected from alkoxy, e.g. OR_{19} or, in the alternative, H, OH, N_3, NHR_{20}, NO_2 or CN, fluoro, chloro, bromo, iodo, where $R_{19\text{-}20}$ are independently selected from the same group which defines $R_{12\text{-}13}$.

14. (original) A compound of claim 3, selected from the group consisting of:





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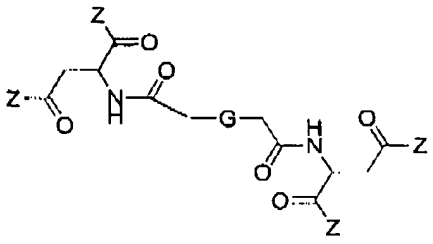
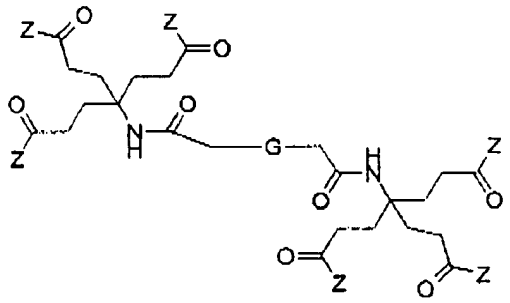
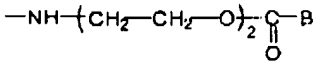


and



wherein Z is one of:



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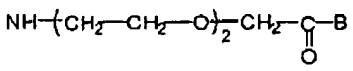

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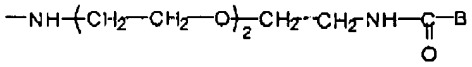

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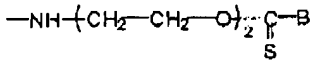

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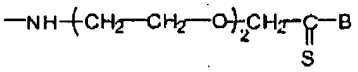

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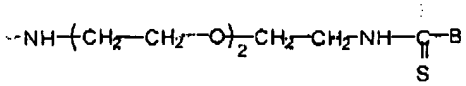

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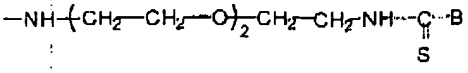
and



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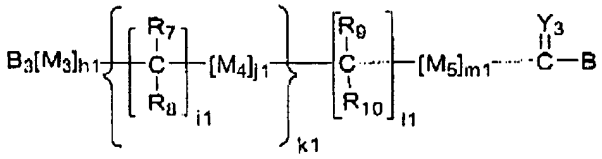
15. (Currently amended) A method of preparing a polymeric transport system, comprising

a) reacting compound of the formula:



wherein

B is a residue of a biologically active amine-containing moiety or a hydroxyl-containing moiety;

B_3 is a cleavable protecting group;

Y_3 is O, S, or NR_{11a};

M_3 and M_4 are independently O, S, or NR_{11b};

M_5 is X or Q;

wherein X is an electron withdrawing group and Q is a moiety containing a free electron pair positioned three to six atoms from $C(=Y_3)$;

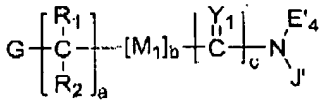
$R_{7\text{-}10}$ and $R_{11a\text{-}b}$ are independently selected from the group consisting of hydrogen, $C_{1\text{-}6}$ alkyls, $C_{3\text{-}12}$ branched alkyls, $C_{3\text{-}8}$ cycloalkyls, $C_{1\text{-}6}$ substituted alkyls, $C_{3\text{-}8}$ substituted cycloalkyls, aryls, substituted aryls, aralkyls, $C_{1\text{-}6}$ heteroalkyls and substituted $C_{1\text{-}6}$ heteroalkyls;

~~*k1-m1*~~ *h1, i1, j1, l1* and *m1* are each independently zero or a positive integer;

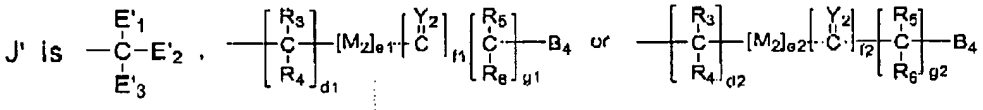
k1 is a positive integer;

b) cleaving the cleavable protecting group B_3; and

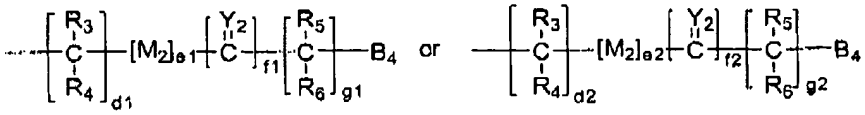
c) reacting the resultant compound with a compound of the formula



wherein

J' is 

E'_{1-4} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} heteroalkyls, C_{1-6} alkoxy, phenoxy, C_{1-6} heteroalkoxy,



wherein

B_4 is a leaving group;

G is a polymer residue;

Y_{1-2} are independently O, S, or NR_{11a};

M_{1-2} are independently O, S, or NR_{11b};

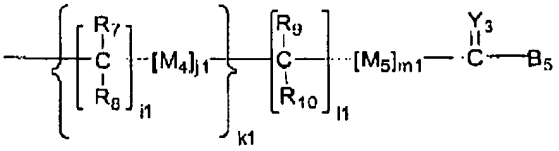
R_{1-6}, R_9 and R_{10} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls and substituted C_{1-6} heteroalkyls;

a, *b*, *c*, d_1-g_1 and d_2-g_2 are each independently zero or a positive integer,

whereby a polymeric conjugate is formed.

16. (Currently amended) A method of preparing a polymeric transport system, comprising:

reacting a biologically active moiety containing an unprotected amino or hydroxyl group with polymeric residue containing a terminal moiety of the formula:



wherein:

Y_3 is O, S, or NR_{11a};

R_{7-10} ~~and NR_{11a}~~ are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls and substituted C_{1-6} heteroalkyls;

M_{4-5} are independently O, S, or NR_{11b};

R_{11a} and R_{11b} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} heteroalkyls, C_{1-6} alkoxy, phenoxy and C_{1-6} heteroalkoxy;

B_5 is a leaving group capable of reacting with an unprotected amino or hydroxyl group of a biologically active moiety; ~~and~~

~~*il ml*~~ *ij, jl, ll* and *ml* are each independently zero or a positive integer, and

kl is a positive integer;

whereby a polymeric conjugate is formed.

17. (original) A method of treatment, comprising:

administering to a mammal in need of such treatment an effective amount of a compound of claim 1, wherein B is a residue of a biologically active moiety.

18. (original) A method of treatment, comprising:

administering to a mammal in need of such treatment an effective amount of a compound of claim 3, wherein B is a residue of a biologically active moiety.